

FILE 'REGISTRY' ENTERED AT 13:26:47 ON 27 MAY 2008

L1 STRUCTURE UPLOADED

L2 11 S L1

L3 157 S L1 SSS FULL

FILE 'CAPLUS' ENTERED AT 13:27:32 ON 27 MAY 2008

L4 2 S L3

=> file registry
COST IN U.S. DOLLARS
FULL ESTIMATED COST

SINCE FILE ENTRY	TOTAL SESSION
0.21	0.21

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STRUCTURE FILE UPDATES: 26 MAY 2008 HIGHEST RN 1022798-85-1
DICTIONARY FILE UPDATES: 26 MAY 2008 HIGHEST RN 1022798-85-1

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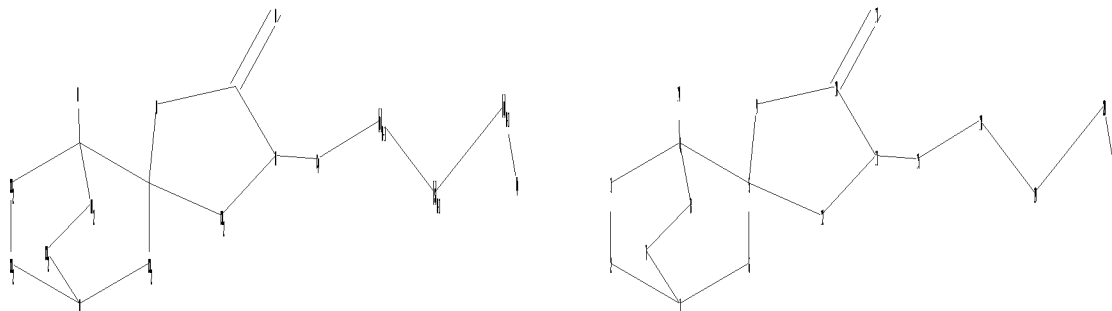
TSCA INFORMATION NOW CURRENT THROUGH January 9, 2008.

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and
predicted properties as well as tags indicating availability of
experimental property data in the original document. For information
on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stndoc/properties.html>

=>
Uploading C:\Program Files\Stnexp\Queries\10563271verify.str



chain nodes :
13 14 15 17 18 19 20
ring nodes :
1 2 3 4 5 6 7 8 9 10 11 12
chain bonds :
4-14 10-13 11-15 15-18 17-20 18-19 19-20
ring bonds :
1-2 1-6 1-7 2-3 3-4 4-5 4-8 5-6 5-9 5-12 7-8 9-10 10-11 11-12
exact/norm bonds :
1-2 1-6 1-7 2-3 3-4 4-5 4-8 5-6 5-9 5-12 7-8 9-10 10-11 10-13 11-12
11-15 15-18 17-20 18-19 19-20
exact bonds :
4-14

G1:O,S,N

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom
11:Atom 12:Atom 13:CLASS 14:CLASS 15:Atom 17:Atom 18:CLASS 19:CLASS
20:CLASS

Generic attributes :

15:

Saturation : Unsaturated

Number of Carbon Atoms : less than 7

Type of Ring System : Monocyclic

17:

Saturation : Unsaturated

Number of Carbon Atoms : less than 7

Type of Ring System : Monocyclic

L1 STRUCTURE UPLOADED

=> s l1

SAMPLE SEARCH INITIATED 13:27:09 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 37 TO ITERATE

100.0% PROCESSED 37 ITERATIONS

11 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 376 TO 1104

PROJECTED ANSWERS: 22 TO 418

L2 11 SEA SSS SAM L1

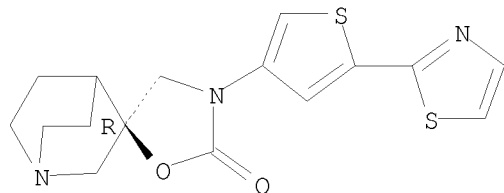
=> d l2 scan

L2 11 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN

IN Spiro[1-azabicyclo[2.2.2]octane-3,5'-oxazolidin]-2'-one,
3'-[5-(2-thiazolyl)-3-thienyl]-, (3R)-

MF C16 H17 N3 O2 S2

Absolute stereochemistry.



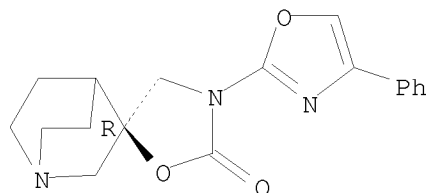
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L2 11 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN

IN Spiro[1-azabicyclo[2.2.2]octane-3,5'-oxazolidin]-2'-one,
3'-(4-phenyl-2-oxazolyl)-, (3R)-
MF C18 H19 N3 O3

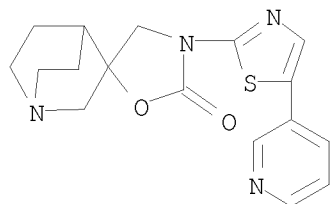
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L2 11 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Spiro[1-azabicyclo[2.2.2]octane-3,5'-oxazolidin]-2'-one,
3'-[5-(3-pyridinyl)-2-thiazolyl]-
MF C17 H18 N4 O2 S



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=> s l1 sss full
FULL SEARCH INITIATED 13:27:26 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 692 TO ITERATE

100.0% PROCESSED 692 ITERATIONS 157 ANSWERS
SEARCH TIME: 00.00.01

L3 157 SEA SSS FUL L1

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	178.36	178.57

FILE 'CAPLUS' ENTERED AT 13:27:32 ON 27 MAY 2008

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FILE COVERS 1907 - 27 May 2008 VOL 148 ISS 22
FILE LAST UPDATED: 26 May 2008 (20080526/ED)

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=> a l3

A IS NOT A RECOGNIZED COMMAND

The previous command name entered was not recognized by the system. For a list of commands available to you in the current file, enter "HELP COMMANDS" at an arrow prompt (=>).

=> s l3

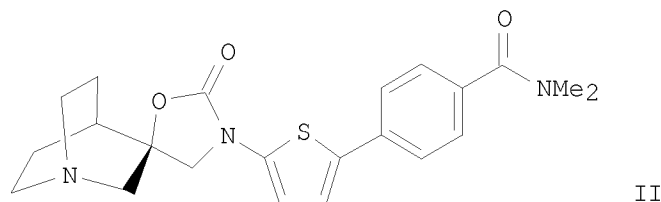
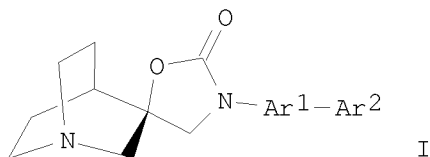
L4 2 L3

=> d l4 ti abs bib

L4 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2008 ACS on STN

TI Preparation of spiro-oxazolidinone compounds as nicotinic acetylcholine receptor ligands

GI



AB Title compds. I [Ar1, Ar2 = 5- or 6-membered aromatic or heteroarom. moiety having 0,1 or 2 nitrogen atoms, 0 or 1 oxygen atoms, and 0 or 1 sulfur atoms; wherein Ar1 is unsubstituted or has 1, 2 or 3 substituents selected from alkyl, alkenyl, alkynyl, etc. and Ar2 is unsubstituted or has 1, 2 or 3 substituents selected from -CONR1R2, -NR1COR2; R1, R2 = H, alkyl, or -NR1R2 in combination is -(CH2)jG(CH2)k-; G = bond, oxygen, sulfur, etc.; j = 2-4; k = 0-2] or stereoisomers, enantiomers, in vivo hydrolysable precursors and pharmaceutically acceptable salts thereof were prepared For example, Pd(PPh3)4 catalyzed coupling reaction of 4-(N,N-dimethylaminocarbonyl)phenylboronic acid with 2,5-dibromothiophene followed by reaction with (3S)-spiro[1-azabicyclo[2.2.2]octane-3,5'-oxazolidin]-2'-one afforded compound II. Compds. I are claimed useful as nicotinic acetylcholine receptor ligands for the treatment of anxiety, schizophrenia, etc. (no data).

AN 2006:608651 CAPLUS <<LOGINID::20080527>>

DN 145:83311

TI Preparation of spiro-oxazolidinone compounds as nicotinic acetylcholine receptor ligands

IN Chapdelaine, Marc; Chang, Hui-Fang; Herzog, Keith J.; Horchler, Carey; Phillips, Eifion

PA Astrazeneca AB, Swed.

SO PCT Int. Appl., 44 pp.

CODEN: PIXXD2

DT Patent

LA English

FAN.CNT 1

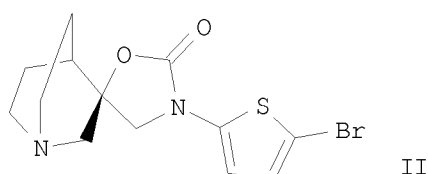
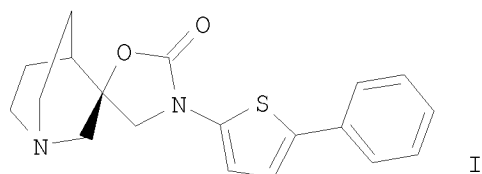
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PI	WO 2006065209	A1	20060622	WO 2005-SE1909	20051213
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	RW: AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LT, LU, LV, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG, BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
	AU 2005317280	A1	20060622	AU 2005-317280	20051213
	CA 2591430	A1	20060622	CA 2005-2591430	20051213
	EP 1831231	A1	20070912	EP 2005-819091	20051213
	R: AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LI, LT, LU, LV, MC, NL, PL, PT, RO, SE, SI, SK, TR, HR				
	MX 200706743	A	20070709	MX 2007-6743	20070606
	IN 2007DN04472	A	20070831	IN 2007-DN4472	20070612
	US 20080113983	A1	20080515	US 2007-721481	20070612
	KR 2007090922	A	20070906	KR 2007-713375	20070614
	NO 2007003551	A	20070801	NO 2007-3551	20070709
	CN 101124232	A	20080213	CN 2005-80048394	20070815
PRAI	US 2004-636362P	P	20041215		
	US 2005-643319P	P	20050112		
	WO 2005-SE1909	W	20051213		

OS MARPAT 145:83311

RE.CNT 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> d 14 2 ti abs bib

L4 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2008 ACS on STN
 TI A preparation of derivatives of oxazolidinone with affinity to the
 $\alpha 7$ -nicotinic acetylcholine receptor
 GI



AB The invention relates to a preparation of derivs. of oxazolidinone of formula Q-X-A-Y [wherein: Q is spiro(azabicyclooctanoxazolidinone) derivative; A is O, S, or NH, etc.; X is 5- or 6-membered heterocycle; Y is 5- or 6-membered (hetero)aromatic ring] with affinity to the $\alpha 7$ -nicotinic acetylcholine receptor. For instance, oxazolidinone derivative I was prepared via phenylation of II by phenylboronic acid. The compds. of the invention were screened in $\alpha 7$ nAChR subtype affinity assay and showed binding affinities (K_i) of less than 1000 nM.

AN 2005:58211 CAPLUS <<LOGINID::20080527>>

DN 142:155977

TI A preparation of derivatives of oxazolidinone with affinity to the $\alpha 7$ -nicotinic acetylcholine receptor

IN Chang, Hui-Fang; Phillips, Eifion

PA Astrazeneca AB, Swed.; Astrazeneca UK Limited

SO PCT Int. Appl., 77 pp.

CODEN: PIXXD2

DT Patent

LA English

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2005005435	A1	20050120	WO 2004-GB2904	20040706
	W:				
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	AU 2004255920	A1	20050120	AU 2004-255920	20040706

CA 2531510	A1	20050120	CA 2004-2531510	20040706
EP 1654264	A1	20060510	EP 2004-743249	20040706
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CN 1829721	A	20060906	CN 2004-80021849	20040706
BR 2004012382	A	20060919	BR 2004-12382	20040706
JP 2007516200	T	20070621	JP 2006-518343	20040706
US 20060154945	A1	20060713	US 2006-563271	20060104
MX 2006PA00231	A	20060411	MX 2006-PA231	20060105
NO 2006000612	A	20060406	NO 2006-612	20060208
PRAI US 2003-485523P	P	20030708		
WO 2004-GB2904	W	20040706		
OS CASREACT 142:155977; MARPAT 142:155977				
RE.CNT 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD				
ALL CITATIONS AVAILABLE IN THE RE FORMAT				